

Iterative Methods for Solving Systems of Linear Equations

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1 – Classic Iterative Methods

1.1 – Basic Concept

First of all we give an example to illustrate the process of iterative methods for solving systems of linear equations.

Consider solving

$$\begin{bmatrix} 3 & 2 \\ 1 & 4 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} 5 \\ 5 \end{bmatrix}.$$

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This system has the exact solution $x_1 = x_2 = 1$. Equivalently we can write the system as

$$\begin{cases} 3x_1 + 2x_2 = 5 \\ x_1 + 4x_2 = 5 \end{cases}$$

This implies that

$$\begin{cases} x_1 = \frac{1}{3}(5 - 2x_2) \\ x_2 = \frac{1}{4}(5 - x_1) \end{cases}$$

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A naive idea is to solve the system by

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that is, to use the iterative formulation

$$\begin{bmatrix} x_1^{(k)} \\ x_2^{(k)} \end{bmatrix} = \begin{bmatrix} \frac{1}{3} & 0 \\ 0 & \frac{1}{4} \end{bmatrix} \left(\begin{bmatrix} 5 \\ 5 \end{bmatrix} - \begin{bmatrix} 0 & 2 \\ 1 & 0 \end{bmatrix} \begin{bmatrix} x_1^{(k-1)} \\ x_2^{(k-1)} \end{bmatrix} \right)$$

If we choose the initial guess $x_1^{(0)} = x_2^{(0)} = 0$, we would obtain

$$\begin{bmatrix} x_1^{(1)} \\ x_2^{(1)} \end{bmatrix} = \begin{bmatrix} \frac{1}{3} & 0 \\ 0 & \frac{1}{4} \end{bmatrix} \left(\begin{bmatrix} 5 \\ 5 \end{bmatrix} - \begin{bmatrix} 0 & 2 \\ 1 & 0 \end{bmatrix} \begin{bmatrix} 0 \\ 0 \end{bmatrix} \right) = \begin{bmatrix} 1.6667 \\ 1.2500 \end{bmatrix}$$

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and

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By repeating the process, we have the following table

k	3	4	5	6	7
$x_1^{(k)}$	1.1111	0.9722	1.0185	0.9954	1.0031
$x_2^{(k)}$	1.0417	0.9722	1.0000	0.9954	1.0012

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This suggests an **iterative process**

$$x^{(k)} = (I - M^{-1}A)x^{(k-1)} + M^{-1}b \equiv Tx^{(k-1)} + c,$$

where T is usually called the **iteration matrix**. The initial vector $x^{(0)}$ can be arbitrary or be chosen according to certain conditions.

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- ☞ $x^{(k)}$ is easily computed. More precisely, the system $Mx^{(k)} = y$ is easy to solve;
- ☞ the sequence $\{x^{(k)}\}$ converges rapidly to the exact solution.

1.2 – Richard's Method

When we choose $M = I$ such that $A = I - (I - A)$, we obtain the iteration procedure

$$x^{(k)} = (I - A)x^{(k-1)} + b = x^{(k-1)} - Ax^{(k-1)} + b \equiv x^{(k-1)} + r^{(k-1)}.$$

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Algorithm 1 (Richard's Method)

for $k = 1, 2, \dots$ **do**

for $i = 1, 2, \dots, n$ **do**

$$r_i^{(k-1)} = b_i - \sum_{j=1}^n a_{ij}x_j^{(k-1)}$$

$$x_i^{(k)} = x_i^{(k-1)} + r_i^{(k-1)}$$

end for

end for

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If we decompose the coefficient matrix A as

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With this method, the iteration matrix $T = -D^{-1}(L + U)$ and $c = D^{-1}b$. Each component $x_i^{(k)}$ can be computed by

$$x_i^{(k)} = \left(b_i - \sum_{j=1}^{i-1} a_{ij}x_j^{(k-1)} - \sum_{j=i+1}^n a_{ij}x_j^{(k-1)} \right) / a_{ii}.$$

$$\begin{aligned} a_{11}x_1^{(k)} + a_{12}x_2^{(k-1)} + a_{13}x_3^{(k-1)} + \cdots + a_{1n}x_n^{(k-1)} &= b_1 \\ a_{21}x_1^{(k-1)} + a_{22}x_2^{(k)} + a_{23}x_3^{(k-1)} + \cdots + a_{2n}x_n^{(k-1)} &= b_2 \\ &\vdots \\ a_{n1}x_1^{(k-1)} + a_{n2}x_2^{(k-1)} + a_{n3}x_3^{(k-1)} + \cdots + a_{nn}x_n^{(k)} &= b_n. \end{aligned}$$

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for $k = 1, 2, \dots$ **do**

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Only the components of $x^{(k-1)}$ are used to compute $x^{(k)}$.

$\Rightarrow x_i^{(k)}, i = 1, \dots, n$, can be computed in **parallel** at each iteration k .

1.4 – Gauss-Seidel Method

When computing $x_i^{(k)}$ for $i > 1$, $x_1^{(k)}, \dots, x_{i-1}^{(k)}$ have already been computed and are likely to be better approximations to the exact x_1, \dots, x_{i-1} than $x_1^{(k-1)}, \dots, x_{i-1}^{(k-1)}$.

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This improvement induce the Gauss-Seidel method.

The Gauss-Seidel method sets $M = D + L$ and defines the iteration as

$$x^{(k)} = -(D + L)^{-1}Ux^{(k-1)} + (D + L)^{-1}b.$$

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Hence each component $x_i^{(k)}$ can be computed by

$$x_i^{(k)} = \left(b_i - \sum_{j=1}^{i-1} a_{ij}x_j^{(k)} - \sum_{j=i+1}^n a_{ij}x_j^{(k-1)} \right) / a_{ii}.$$

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Algorithm 3 (Gauss-Seidel Method)

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Hence the iteration matrix $T = (D + \omega L)^{-1}[(1 - \omega)D - \omega U]$. Each component $x_i^{(k)}$ can be computed by the formulation

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The question of choosing a good relaxation parameter ω is a very complex topic.

1.6 – Symmetric Successive Over Relaxation (SSOR) Method

In theory the symmetric successive over relaxation (SSOR) method chooses the splitting matrix $M = \frac{1}{\omega(2-\omega)}(D + \omega L)D^{-1}(D + \omega U)$ and iterates with the iteration matrix

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The idea is in fact to implement the SOR formulation **twice**, **one forward** and **one backward**, at each iteration. That is, SSOR method defines

$$\begin{aligned}(D + \omega L)x^{(k-\frac{1}{2})} &= ((1 - \omega)D - \omega U)x^{(k-1)} + \omega b \\(D + \omega U)x^{(k)} &= ((1 - \omega)D - \omega L)x^{(k-\frac{1}{2})} + \omega b\end{aligned}$$

Each component $x_i^{(k)}$ is obtained by first computing

$$x_i^{(k-\frac{1}{2})} = \omega \left(b_i - \sum_{j=1}^{i-1} a_{ij} x_j^{(k-\frac{1}{2})} - \sum_{j=i+1}^n a_{ij} x_j^{(k-1)} \right) / a_{ii} + (1 - \omega) x_i^{(k)}$$

followed by

$$x_i^{(k)} = \omega \left(b_i - \sum_{j=1}^{i-1} a_{ij} x_j^{(k)} - \sum_{j=i+1}^n a_{ij} x_j^{(k-\frac{1}{2})} \right) / a_{ii} + (1 - \omega) x_i^{(k-\frac{1}{2})}.$$

2 – Convergence Analysis

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$$\rho(A) = \max |\lambda| \leq \|A\|. \quad \blacksquare$$

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Next we show that $(I - A)^{-1} = I + A + A^2 + \dots$.

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Lemma 3 Suppose that $A \in \mathbb{R}^{n \times n}$ and $\|\cdot\|$ is a subordinate matrix norm. If $\|A\| < 1$, then $I - A$ is nonsingular and

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This shows that $(I - A)^{-1} = \sum_{k=0}^{\infty} A^k$.

Finally, since $\|A\| < 1$,

$$\|(I - A)^{-1}\| = \left\| \sum_{k=0}^{\infty} A^k \right\| \leq \sum_{k=0}^{\infty} \|A^k\| \leq \sum_{k=0}^{\infty} \|A\|^k = \frac{1}{1 - \|A\|}.$$



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Theorem 2 *The following statements are equivalent.*

1. A is a *convergent matrix*, i.e., $A^k \rightarrow 0$ as $k \rightarrow \infty$;
2. $\lim_{k \rightarrow \infty} \|A^k\| = 0$ for *some* subordinate matrix norm;
3. $\lim_{k \rightarrow \infty} \|A^k\| = 0$ for *all* subordinate matrix norm;
4. $\rho(A) < 1$;
5. $\lim_{k \rightarrow \infty} A^k x = 0$ for *any* x .

Theorem 3 For *any* $x^{(0)} \in \mathbb{R}^n$, the sequence produced by

$$x^{(k)} = Tx^{(k-1)} + c, \quad k = 1, 2, \dots,$$

converges to the *unique* solution of $x = Tx + c$ if and only if

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$$x^{(1)} = Tx^{(0)} + c$$

$$x^{(2)} = Tx^{(1)} + c = T^2x^{(0)} + (T + I)c$$

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In general

$$x^{(k)} = T^k x^{(0)} + (T^{k-1} + T^{k-2} + \dots + T + I)c.$$

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It follows from theorem $\rho(T) < 1$. ■

Corollary 1 *If $\|T\| < 1$ for **some** subordinate matrix norm, then the sequence produced by*

$$x^{(k)} = Tx^{(k-1)} + c$$

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Proof: Since $x = Tx + c$ and $x^{(k)} = Tx^{(k-1)} + c$,

$$\begin{aligned}x - x^{(k)} &= Tx + c - Tx^{(k-1)} - c \\&= T(x - x^{(k-1)}) \\&= T^2(x - x^{(k-2)}) = \dots = T^k(x - x^{(0)}).\end{aligned}$$

The first statement can then be derived

$$\|x - x^{(k)}\| = \|T^k(x - x^{(0)})\| \leq \|T\|^k \|x - x^{(0)}\|.$$

For the second result, we first show that

$$\|x^{(n)} - x^{(n-1)}\| \leq \|T\|^{n-1} \|x^{(1)} - x^{(0)}\| \text{ for any } n \geq 1.$$

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Let $m \geq k$,

$$\begin{aligned} &x^{(m)} - x^{(k)} \\ &= \left(x^{(m)} - x^{(m-1)}\right) + \left(x^{(m-1)} - x^{(m-2)}\right) + \dots + \left(x^{(k+1)} - x^{(k)}\right) \\ &= T^{m-1} \left(x^{(1)} - x^{(0)}\right) + T^{m-2} \left(x^{(1)} - x^{(0)}\right) + \dots + T^k \left(x^{(1)} - x^{(0)}\right) \\ &= \left(T^{m-1} + T^{m-2} + \dots + T^k\right) \left(x^{(1)} - x^{(0)}\right), \end{aligned}$$

hence

$$\begin{aligned} & \|x^{(m)} - x^{(k)}\| \\ \leq & (\|T\|^{m-1} + \|T\|^{m-2} + \cdots + \|T\|^k) \|x^{(1)} - x^{(0)}\| \end{aligned}$$

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This proves the second result. ■

Theorem 6 If A is *strictly diagonal dominant*, then both the *Jacobi* and *Gauss-Seidel* methods *converges* for *any* initial vector $x^{(0)}$.

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For Jacobi method, the iteration matrix $T_J = -D^{-1}(L + U)$ has entries

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Hence

$$\|T_J\|_{\infty} = \max_{1 \leq i \leq n} \sum_{j=1, j \neq i}^n \left| \frac{a_{ij}}{a_{ii}} \right| = \max_{1 \leq i \leq n} \frac{1}{|a_{ii}|} \sum_{j=1, j \neq i}^n |a_{ij}| < 1,$$

and this implies that the Jacobi method converges.

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Hence for $i = 1, \dots, n$,

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Choose the index k such that $|y_k| = 1 \geq |y_j|$ (this index can always be found since $\|y\|_\infty = 1$).

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Since λ is arbitrary, $\rho(T_{GS}) < 1$.

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Since λ is arbitrary, $\rho(T_{GS}) < 1$. This means the Gauss-Seidel method converges. ■

Theorem 7 If A is *positive definite* and the relaxation parameter ω satisfying $0 < \omega < 2$, then the *SOR* iteration *converges* for *any* initial vector $x^{(0)}$.

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Theorem 8 If A is *positive definite* and *tridiagonal*, then $\rho(T_{GS}) = [\rho(T_J)]^2 < 1$ and the *optimal* choice of ω for the *SOR* iteration is

$$\omega = \frac{2}{1 + \sqrt{1 - [\rho(T_J)]^2}}.$$

With this choice of ω , $\rho(T_{SOR}) = \omega - 1$.